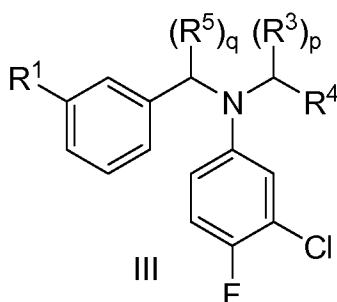


IN THE CLAIMS:

1. Cancelled
2. Cancelled
3. (Currently amended) The compound according to Claim 2 of the Formula III;



wherein:

R<sup>1</sup> is selected from: H, F and OH;

R<sup>3</sup> is selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C<sub>2</sub>-C<sub>6</sub>)alkenyl and (C<sub>1</sub>-C<sub>6</sub>)alkyl hydroxyl;

R<sup>4</sup> is selected from: H, oxo, OH, halo, CN, NH<sub>2</sub>, NO<sub>2</sub>, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>2</sub>-C<sub>10</sub>)alkenyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>2</sub>-C<sub>10</sub>)alkynyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>, C(O)H, (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H, C(O)N(R<sup>b</sup>)<sub>2</sub>, and S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>; said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>;

R<sup>5</sup> is H or CH<sub>3</sub>;

R<sup>a</sup> is selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl and heterocyclyl; said alkyl, cycloalkyl, aryl and heterocyclyl is optionally substituted with one or more substituents selected from OH, (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, (O)C=O(C<sub>1</sub>-C<sub>6</sub>)alkyl, oxo and N(R<sup>c</sup>)<sub>2</sub>;

R<sup>b</sup> is independently selected from: H, oxo, OH, halogen, CO<sub>2</sub>H, CN, (O)C=O(C<sub>1</sub>-C<sub>6</sub>)alkyl, N(R<sup>c</sup>)<sub>2</sub>, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)O(C<sub>1</sub>-C<sub>6</sub>)alkyl, C=O(C<sub>1</sub>-C<sub>6</sub>)alkyl and S(O)<sub>2</sub>R<sup>a</sup>; said alkyl, cycloalkyl, aryl or heterocyclyl is optionally substituted with one or more substituents selected from OH, (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, (O)C=O(C<sub>1</sub>-C<sub>6</sub>)alkyl, oxo, N(R<sup>c</sup>)<sub>2</sub> and optionally substituted heterocyclyl, wherein said heterocyclyl is optionally substituted with (C<sub>1</sub>-C<sub>6</sub>)alkyl, oxo or NH<sub>2</sub>.

R<sup>c</sup> is independently selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl;

a is 0 or 1;

b is 0 or 1;

p is 1 or 2;

q is 0 or 1;

~~and all other substituents and variables are as defined in Claim 2;~~

or a pharmaceutically acceptable salt or stereoisomer thereof.

4. (Original) The compound according to Claim 3 of the Formula III;

wherein:

R<sup>4</sup> is selected from: H, oxo, OH, halo, CN, NH<sub>2</sub>, NO<sub>2</sub>, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>2</sub>-C<sub>10</sub>)alkenyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>2</sub>-C<sub>10</sub>)alkynyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>, C(O)H, (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H, C(O)N(R<sup>b</sup>)<sub>2</sub>, and S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>; said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>;

and all other substituents and variables are as defined in Claim 3;

or a pharmaceutically acceptable salt or stereoisomer thereof.

5. (Original) The compound according to Claim 3 of the Formula III;

wherein:

R<sup>4</sup> is selected from: H, oxo, OH, halo, CN, NH<sub>2</sub>, NO<sub>2</sub>, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>2</sub>-C<sub>10</sub>)alkenyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>2</sub>-C<sub>10</sub>)alkynyl, (C=O)<sub>a</sub>O<sub>b</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>, (C=O)-R<sup>b</sup>, C(O)H, (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H, C(O)N(R<sup>b</sup>)<sub>2</sub>, and S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>; said alkyl, alkenyl, alkynyl and alkylene is optionally substituted with up to three substituents selected from R<sup>b</sup>;

and all other substituents and variables are as defined in Claim 3;

or a pharmaceutically acceptable salt or stereoisomer thereof.

6. (Original) A compound which is selected from:

3-[(3-chloro-4-fluorophenyl)(2-hydroxy-1-methylethyl)amino]methyl} phenol;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]propan-1-ol;  
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)-N-1-[2(dimethylamino)ethyl] alaninamide;  
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)alaninamide;  
methyl N-benzyl-N-(3-chloro-4-fluorophenyl)alanylglycinate;  
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)-N-1-(isoxazol-4-ylmethyl)alaninamide;  
3-[benzyl(3-chloro-4-fluorophenyl)amino]-2-methylbutan-2-ol;  
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)-N-1,N-1-dimethylpropane-1,2-diamine;  
N-benzyl-3-chloro-4-fluoro-N-[1-methyl-2-(4-methylpiperazin-1-yl)ethyl]aniline;  
2-[(3-chloro-4-fluorophenyl)(1-phenylethyl)amino]propan-1-ol;  
N-2-(3-chloro-4-fluorophenyl)-N-2-(3-hydroxybenzyl)-N-1-(isoxazol-4-ylmethyl)alaninamide;  
N-2-(3-chloro-4-fluorophenyl)-N-2-(3-hydroxybenzyl)alaninamide;  
N-2-(3-chloro-4-fluorophenyl)-N-1-[2-(dimethylamino)ethyl]-N-2-(3-hydroxybenzyl)alaninamide;  
Methyl 2-[benzyl(3-chloro-4-fluorophenyl)amino]butanoate;  
Methyl 2-[benzyl(3-chloro-4-fluorophenyl)amino]pent-4-enoate;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]pent-4-en-1-ol;  
N-benzyl-N-(3-chloro-4-fluorophenyl)glycine;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]pentan-1-ol;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]butan-1-ol;  
N-benzyl-3-chloro-N-[1-(3-[(dimethylamino)methyl]piperidin-1-yl)carbonyl]propyl]-4-fluoroaniline;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]-N-methyl-N-[2-(1-methyl-1H-pyrazol-4-yl)ethyl]butanamide;

2-[benzyl(3-chloro-4-fluorophenyl)amino]-3-methylbutan-1-ol;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]pentane-1,5-diol;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]-3-cyclopropylpropan-1-ol; and  
N<sup>2</sup>-benzyl-N<sup>2</sup>-(3-chloro-4-fluorophenyl)-N<sup>1</sup>-[2-(dimethylamino)ethyl]-2-methylalaninamide;

or a pharmaceutically acceptable salt or stereoisomer thereof.

7. (Currently amended) The TFA salt of a compound according to Claim 4 3 which is

3-[(3-chloro-4-fluorophenyl)(2-hydroxy-1-methylethyl)amino]methyl} phenol;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]propan-1-ol;  
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)-N-1-[2(dimethylamino)ethyl] alaninamide;  
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)alaninamide;  
methyl N-benzyl-N-(3-chloro-4-fluorophenyl)alanyl glycinate;  
3-[benzyl(3-chloro-4-fluorophenyl)amino]-2-methylbutan-2-ol;  
2-[(3-chloro-4-fluorophenyl)(1-phenylethyl)amino]propan-1-ol;  
N-2-(3-chloro-4-fluorophenyl)-N-2-(3-hydroxybenzyl)-N-1-(isoxazol-4-ylmethyl)alaninamide;  
N-2-(3-chloro-4-fluorophenyl)-N-2-(3-hydroxybenzyl)alaninamide;  
N-2-(3-chloro-4-fluorophenyl)-N-1-[2-(dimethylamino)ethyl]-N-2-(3-hydroxybenzyl)alaninamide;  
N-benzyl-N-(3-chloro-4-fluorophenyl)glycine;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]pentan-1-ol;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]butan-1-ol;  
N-benzyl-3-chloro-N-[1-(3-[(dimethylamino)methyl]piperidin-1-yl} carbonyl)propyl]-4-fluoroaniline;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]-N-methyl-N-[2-(1-methyl-1H-pyrazol-4-yl)ethyl]butanamide;  
2-[benzyl(3-chloro-4-fluorophenyl)amino]pentane-1,5-diol; and  
N<sup>2</sup>-benzyl-N<sup>2</sup>-(3-chloro-4-fluorophenyl)-N<sup>1</sup>-[2-(dimethylamino)ethyl]-2-methylalaninamide;

or stereoisomer thereof.

8. (Currently amended) A pharmaceutical composition comprising a pharmaceutical carrier, and dispersed therein, a therapeutically effective amount of a compound of Claim 34.

9. (Currently amended) ~~The use of the compound according to Claim 1 for the preparation of a medicament useful in the treatment or prevention of~~ A method of treating cancer in a mammal in need of such treatment with a compound of Claim 3.